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B109 Isolation of mangiferin and structure revision of shamimin from Bombax malabaricum

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Bombax malabaricum DC. (Bombacaceae) (Syn. B. ceiba L. and Salmalia malabaricum DC) is also known as silkcotton tree. It is commonly found in the Indo-Pakistan subcontinent and other parts of Asia and Australia (1). The plant is well reputed for the treatment of diarrhoea, tumors, fever, dysentery, kidney and bladder ulceration, and chronic inflammation (2,3). Phytochemical investigation of different parts of this plant resulted in the isolation of naphthol, naphtoquinones, polysaccharides, anthocyanins and lupeol (4).

The 70 % alcoholic extract of the leaves of *B. malabaricum* was concentrated, kept in the refrigerator overnight, and centrifugated. The supernatant was successively extracted with CHCl₃, EtOAc and BuOH. Repeated column chromatography of the BuOH fraction yielded compound **1** which was identified as $2C\beta$ -D-glucosyl-1,3,6,7-tetrahydroxyxanthone (mangiferin) (4,5). The ¹H and ¹³C-NMR data were in complete agreement with those reported for shamimin or $6C\beta$ -D-glucosyl-3,5,7,2',4',5'-hexahydroxyflavone reported before from *Bombax ceiba* (3). Therefore we conclude that the structure of shamimin has to be revised, and that it is identical to mangiferin.

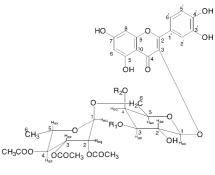
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B110 Structure elucidation of three new acetylated flavonoid glycosides from Centaurium spicatum

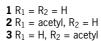
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Centaurium spicatum (L.) Fritsch (Gentianaceae) is an annual herb occurring in Southern Europe and Northern Africa, where it is used together with other Centaurium species like C. pulchellum in traditional medicine. Alkaloids

and secoiridoids have been reported before from Centaurium spicatum, but this is the first report on flavonoids from this plant. Three new acetylated flavonol glycosides, quercetin 30(2,3,4-triacetyl-arhamnopyranosyl)-(1 \rightarrow 6)-β-galacto-pyranoside (1), quercetin 3-O-[(2,3,4-triacetyl-galactopyranosyl)-(1-6)]-3-acetyl-β-galactopyranoside (2), and quercetin 3-O-[(2,3,4-triacetyl- α -rhamnopyranosyl)-(1 \rightarrow 6)]-4-acetyl-β-galactopyranoside (3) have been isolated and identified. Structure elucidation, especially the localisation of the acetyl groups, and complete ¹H and ¹³C NMR assignments were carried out using one- and twodimensional NMR methods, including ¹H and ¹³C NMR, DEPT-135 and DEPT-90, and gradient-assisted experiments such as DQF-COSY, TOCSY, HSQC and HMBC (1).



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50th Annual Congress of the Society for Medicinal Plant Research